

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112320\
 Data File : BM027919.D
 Acq On : 23 Nov 2020 18:37
 Operator : JU/CG
 Sample : L4711-13DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B65DL

Manual Integrations
 APPROVED

mohammad
 11/25/2020 12:44:48 PM

Quant Time: Nov 24 00:52:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 24 00:42:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1137	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	4597	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2795	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	5477	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	3285	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2745	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	425	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	734	0.05	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.22	128	404	0.029	ng/ul#	68
5) 2-Methylnaphthalene	12.80	142	200	0.021	ng/ul	95
7) Acenaphthylene	14.67	152	328	0.024	ng/ul#	60
8) Acenaphthene	15.00	153	729	0.071	ng/ul	97
9) Fluorene	15.98	166	1082	0.091	ng/ul#	98
12) Phenanthrene	17.70	178	43026	2.487	ng/ul	97
13) Anthracene	17.79	178	6182	0.370	ng/ul	100
15) Fluoranthene	19.68	202	117899	6.146	ng/ul	100
17) Pyrene	20.04	202	93079	5.579	ng/ul	99
18) Benzo(a)anthracene	21.75	228	41135	3.104	ng/ul	100
19) Chrysene	21.80	228	47300	3.664	ng/ul	99
21) Benzo(b)fluoranthene	23.45	252	53762m	4.660	ng/ul	
22) Benzo(k)fluoranthene	23.49	252	17866m	1.581	ng/ul	
23) Benzo(a)pyrene	24.08	252	34227	3.393	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.68	276	22588	2.068	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.68	278	5581	0.614	ng/ul	96
26) Benzo(a,h,i)perylene	27.45	276	21509	2.348	ng/ul	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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